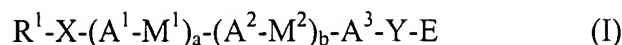


**IN THE CLAIMS:**

1. (Currently Amended) A five-membered ring compound of the formula (I),



where the symbols and indices have the following meanings:

**E** is a radical T-Z-R<sup>2</sup> containing a five-membered ring, where:

(i) **T** is undirected and is  
4-fluorothiophene-2,5-diyl, 3-fluorothiophene-2,5-diyl,  
3-fluorothiophene-2,4-diyl or 5-fluorothiophene-2,4-diyl

**Z** is a single bond or -O-

**R<sup>2</sup>** is hydrogen or a straight-chain or branched alkyl radical (with or without asymmetric carbon atoms) having 1 to 20 carbon atoms, where one nonterminal CH<sub>2</sub> group may be replaced by -O- or -OC(=O)- or -C(=O)O- and/or one or more H atoms may be replaced by F, with the provisos that  
a) the -CH<sub>2</sub>- group nearest to the thiophene cannot be replaced by -O- when Z is -O-  
b) R<sup>2</sup> can only be hydrogen when Z is a single bond,

**Y** is -OC(=O)-, -OCH<sub>2</sub>-, -CH<sub>2</sub>CH<sub>2</sub>-

**a, b** are each, independently of one another, 0 or 1

(ii) **T** is furan-2,5-diyl or furan-2,4-diyl

**Z** is a single bond or -O-

**R<sup>2</sup>** is a straight-chain or branched alkyl radical (with or without asymmetric carbon atoms) having 1 to 20 carbon atoms, where one nonterminal CH<sub>2</sub>

group nonadjacent to furan may be replaced by -O- or -OC(=O)- or -C(=O)O- and/or one or more H atoms may be replaced by F,

Y is -OC(=O)-, -OCH<sub>2</sub>-, -CH<sub>2</sub>CH<sub>2</sub>-

a, b are each, independently of one another, 0 or 1

- (iii) **T** is undirected and is isoxazole-3,5-diyl  
**Z** is a single bond or -O-  
**R<sup>2</sup>** is hydrogen or a straight-chain or branched alkyl radical (with or without asymmetric carbon atoms) having 1 to 20 carbon atoms, where one nonterminal CH<sub>2</sub> group may be replaced by -O- or -OC(=O)- or -C(=O)O- and/or one or more H atoms may be replaced by F, with the provisos that  
a) the -CH<sub>2</sub>- group nearest to the isoxazole cannot be replaced by -O- when Z is -O-  
b) R<sup>2</sup> can only be hydrogen when Z is a single bond,

a is 1

b is 0 or 1

Y is -OC(=O)-, -OCH<sub>2</sub>-, -CH<sub>2</sub>CH<sub>2</sub>-

- (iv) **T** is undirected and is thiazole-2,5-diyl or thiazole-2,4-diyl  
**Z** is a single bond  
**R<sup>2</sup>** is hydrogen or a straight-chain or branched alkyl radical (with or without asymmetric carbon atoms) having 1 to 20 carbon atoms, where one nonterminal CH<sub>2</sub> group may be replaced by -O- or -OC(=O)- or -C(=O)O- and/or one or more H atoms may be replaced by F,

Y is -OC(=O)-, -OCH<sub>2</sub>-, -CH<sub>2</sub>CH<sub>2</sub>-

a, b are each, independently of one another, 0 or 1

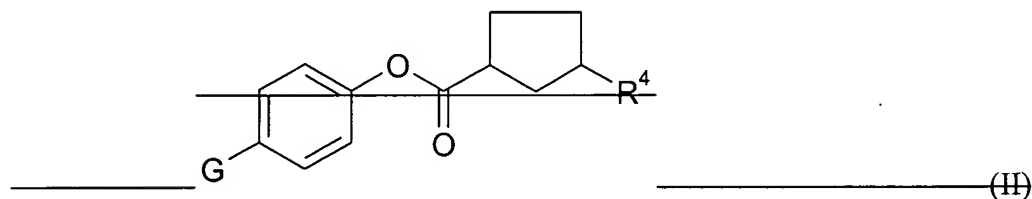
- (v) **T** is cyclopentane-1,3-diyl

~~Z~~ is a single bond or ~~O~~

~~R<sup>2</sup>~~ is hydrogen or a straight chain or branched alkyl radical (with or without asymmetric carbon atoms) having 1 to 20 carbon atoms, where one nonterminal CH<sub>2</sub> group may be replaced by ~~O~~ or ~~OC(=O)~~ or ~~C(=O)O~~ and/or one or more H atoms may be replaced by F, with the provisos that  
a) the ~~CH<sub>2</sub>~~ group nearest to the cyclopentane cannot be replaced by ~~O~~ when Z is ~~O~~

b) ~~R<sup>2</sup>~~ can only be hydrogen when Z is a single bond,

with the exception of compounds of the formula (II)



in which

~~R<sup>4</sup>~~ is as defined for ~~R<sup>2</sup>~~

~~G~~ is trans 4 propyl cyclohexyl or trans 4 butyl cyclohexyl or an alkyl group of 1 to 15 carbon atoms, in which, in addition, one or more nonadjacent CH<sub>2</sub> groups may be replaced by ~~O~~, ~~CO~~, ~~OCO~~, ~~O CO O~~, ~~CHhalogen~~, ~~CHCN~~ and/or ~~CH=CH~~ or is F, CN,

(vi) T is cyclopentane-1,3-diyl, in which one -CH<sub>2</sub>CH<sub>2</sub>- or -CH<sub>2</sub>CH- group is replaced by a -CH=CH- or CH=C- group respectively

Z is a single bond

R<sup>2</sup> is hydrogen or a straight-chain or branched alkyl radical (with or without asymmetric carbon atoms) having 1 to 20 carbon atoms, where one nonterminal CH<sub>2</sub> group may be replaced by -O- or -OC(=O)- or -C(=O)O- and/or one or more H atoms may be replaced by F, with the proviso that the -CH<sub>2</sub>- group nearest to the cyclopentene cannot be replaced and where

Y cannot be -CH<sub>2</sub>-CH<sub>2</sub>-,

a is 1

b is 0 or 1

Y is -OC(=O)-, -OCH<sub>2</sub>-

**R<sup>1</sup>** is hydrogen or a straight-chain or branched C<sub>1-20</sub>-alkyl or C<sub>2-20</sub>-alkenyl radical (with or without asymmetric carbon atoms), where

- a) one or two nonterminal CH<sub>2</sub> groups may be replaced, independently of one another, by -O- or -C(=O)-, with the proviso that two adjacent CH<sub>2</sub> groups cannot be replaced in the same way, and/or
- b) one CH<sub>2</sub> group may be replaced by -C≡C-, and/or
- c) one CH<sub>2</sub> group may be replaced by -Si(CH<sub>3</sub>)<sub>2</sub>-, cyclopropane-1,2-diyl, cyclobutane-1,3-diyl, cyclopentane-1,4-diyl, bicyclo[1.1.1]pentane-1,3-diyl or cyclohexane-1,4-diyl, and/or
- d) one or more H atoms may be replaced by F and/or CN,
- e) in the case of a branched alkyl radical containing asymmetric carbon atoms, the asymmetric carbon atoms have -CH<sub>3</sub>, -OCH<sub>3</sub>, -CF<sub>3</sub>, F, CN and/or Cl as substituents or are incorporated into a 3- to 7-membered ring, in which, in addition, one or two non-adjacent CH<sub>2</sub> groups may be replaced by -O- and one CH<sub>2</sub> group non-adjacent to these groups may be replaced by -OC(=O)-;

**X** is a single bond, -O-, OC(=O)-, -C(=O)O- or -OC(=O)O-

~~**Y** is -OC(=O)-, -OCH<sub>2</sub>-, -CH<sub>2</sub>CH<sub>2</sub>-~~

**A<sup>1</sup>, A<sup>2</sup>, A<sup>3</sup>** are each, independently of one another, phenylene-1,4-diyl, unsubstituted or monosubstituted or disubstituted by CN or F, phenylene-1,3-diyl, unsubstituted or monosubstituted or disubstituted by CN or F, cyclohexane-1,4-diyl, in which one or two H atoms may be replaced by CN and/or CH<sub>3</sub> and/or F, 1-cyclohexene-1,4-diyl, in which one H atom may be replaced by F, 1-alkyl-1-silacyclohexane-1,4-diyl, pyridine-2,5-diyl, unsubstituted or monosubstituted by F, pyrimidine-2,5-

diyl, unsubstituted or monosubstituted by F, cyclopentane-2,5-diyl or thiophene-2,5-diyl;

$M^1$ ,  $M^2$  are undirected and are each, independently of one another, -OC(=O)-, -OCH<sub>2</sub>-, -CH<sub>2</sub>CH<sub>2</sub>-, -OC(=O)CH<sub>2</sub>CH<sub>2</sub>-, -OCH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>-, -C≡C-, -CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>- or a single bond;

~~a, b are each, independently of one another, 0 or 1.~~

2. (Original) A liquid-crystal mixture comprising at least one compound of the formula (I) as claimed in claim 1.

3. (Original) A liquid-crystal mixture as claimed in claim 2, which comprises from 0.01 to 80% by weight of one or more compounds of the formula (I).

4. (Previously presented) A liquid-crystal mixture as claimed in claim 2, which is ferroelectric (chiral smectic).

5. (Previously presented) A liquid-crystal mixture as claimed in claim 2, which is nematic.

6. (Original) A ferroelectric switching and/or display device, which contains a ferroelectric liquid-crystal mixture as claimed in claim 4.

7. (Original) A ferroelectric switching and/or display device as claimed in claim 6, which contains active matrix elements and wherein the liquid-crystal layer forms a monostable monodomain.

8. (Cancelled)

9. (Cancelled)

10. (Cancelled)

11. (Cancelled)

12. (Cancelled)